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spectra
NEWS 4 MAR 31 CA/Caplus and CASREACT patent number format for U.S.
applications updated
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NEWS 6 MAR 31 EMBASE, EMBAL, and LEMBASE reloaded with enhancements
NEWS 7 APR 04 STN AnaVist, Version 1, to be discontinued
NEWS 8 APR 15 WPIDS, WPINDEX, and WPIX enhanced with new
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NEWS 10 APR 28 IMSRESEARCH reloaded with enhancements
NEWS 11 MAY 30 INPAFAMDB now available on STN for patent family
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NEWS 14 JUN 06 KOREAPAT updated with 41,000 documents
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NEWS 21 JUN 30 STN AnaVist enhanced with database content from EPFULL
NEWS 22 JUL 28 CA/Caplus patent coverage enhanced
NEWS 23 JUL 28 EPFULL enhanced with additional legal status
information from the epoline Register
NEWS 24 JUL 28 IFICDB, IFIPAT, and IFIUDB reloaded with enhancements
NEWS 25 JUL 28 STN Viewer performance improved

NEWS EXPRESS JUNE 27 08 CURRENT WINDOWS VERSION IS V8.3,
AND CURRENT DISCOVER FILE IS DATED 23 JUNE 2008.

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COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	0.21	0.21

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STRUCTURE FILE UPDATES: 27 JUL 2008 HIGHEST RN 1036536-16-9
DICTIONARY FILE UPDATES: 27 JUL 2008 HIGHEST RN 1036536-16-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

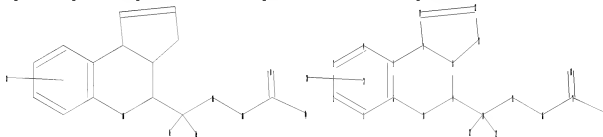
TSCA INFORMATION NOW CURRENT THROUGH January 9, 2008.

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=>
Uploading C:\Program Files\Stnexp\Queries\10-518405genBb.str



chain nodes :
14 15 16 17 18 19 20 22 23
ring nodes :

```

1  2  3  4  5  6  7  8  9  10  11  12  13
chain bonds :
9-14  14-15  14-22  14-23  15-16  16-17  17-18  17-19
ring bonds :
1-2  1-6  2-3  3-4  4-5  5-6  5-7  6-10  7-8  7-11  8-9  8-13  9-10  11-12  12-13
exact/norm bonds :
5-7  6-10  7-8  7-11  8-9  8-13  9-10  11-12  12-13  14-15  15-16  16-17  17-18
17-19
exact bonds :
9-14  14-22  14-23
normalized bonds :
1-2  1-6  2-3  3-4  4-5  5-6

```

Match level :

```

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:Atom 12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS
19:CLASS 20:CLASS 21:Atom 22:CLASS 23:CLASS

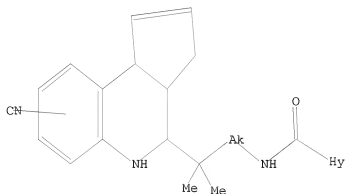
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L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> s l1 sss full

FULL SEARCH INITIATED 17:46:15 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 621 TO ITERATE

100.0% PROCESSED 621 ITERATIONS

13 ANSWERS

SEARCH TIME: 00.00.01

L2 13 SEA SSS FUL L1

=> file caplus

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

178.36

178.57

FILE 'CAPLUS' ENTERED AT 17:46:25 ON 28 JUL 2008

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FILE COVERS 1907 - 28 Jul 2008 VOL 149 ISS 5
FILE LAST UPDATED: 27 Jul 2008 (20080727/ED)

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<http://www.cas.org/legal/infopolicy.html>

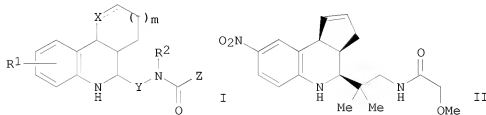
=> s l2

L3 1 L2

=> d l3 abs ibib hitstr

L3 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2008 ACS on STN

GI



AB The title nonsteroidal tetrahydroquinoline derivs. with general formula of I [wherein R1 = NO2 or CN; X = CH or O; m = 0 or 1; Y = (un)substituted alkylene; R2 = H, alkyl, cycloalkyl, or aralkyl; Z = (un)substituted alkyl, aryl, etc.] or pharmaceutically acceptable salts thereof are prepared as androgen receptor agonists. For example, the compound II was prepared in a three-step synthesis starting from 4-nitroaniline, cyclopentadiene, and tert-Bu N-(2,2-dimethyl-3-oxopropyl)carbamate. II showed relative binding affinity of 1076 against androgen receptor in rat. Formulations containing I as an active ingredient were also described.

ACCESSION NUMBER: 2004:2862 CAPLUS

DOCUMENT NUMBER: 140:59527

TITLE: Preparation of bicyclic tetrahydroquinoline

derivatives as androgen receptor agonists

INVENTOR(S): Miyakawa, Motonori; Sumita, Yuji; Furuya, Kazuyuki;
Ichikawa, Kiyonoshin; Yamamoto, Noriko; Hanada, Keigo;
Amano, Seiji; Nejishima, Hiroaki

PATENT ASSIGNEE(S): Kaken Pharmaceutical Co., Ltd., Japan
 SOURCE: PCT Int. Appl., 85 pp.
 CODEN: PIXXD2
 DOCUMENT TYPE: Patent
 LANGUAGE: Japanese
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004000816	A1	20031231	WO 2003-JP7799	20030619
W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW			
RW:	GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
AU 2003244313	A1	20040106	AU 2003-244313	20030619
EP 1520856	A1	20050406	EP 2003-760911	20030619
R:	AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK			
US 20060128737	A1	20060615	US 2005-518405	20051118
PRIORITY APPLN. INFO.:			JP 2002-179088	A 20020619
			WO 2003-JP7799	W 20030619

OTHER SOURCE(S): MARPAT 140:59527

IT 637333-89-2P 637333-90-5P 637333-91-6P
 637333-92-7P 637333-93-8P 637333-94-9P
 637333-95-0P 637333-96-1P 637333-97-2P
 637333-98-3P 637333-99-4P 637334-00-0P
 637334-01-1P

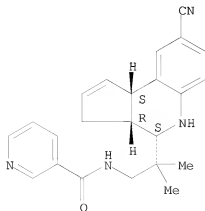
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(drug candidate; preparation of bicyclic tetrahydroquinoline derivs. as androgen receptor agonists)

RN 637333-89-2 CAPLUS

CN 3-Pyridinecarboxamide, N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-, rel- (CA INDEX NAME)

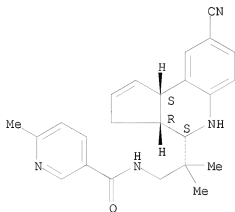
Relative stereochemistry.



RN 637333-90-5 CAPLUS

CN 3-Pyridinecarboxamide, N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-6-methyl-, rel- (CA INDEX NAME)

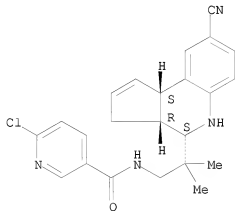
Relative stereochemistry.



RN 637333-91-6 CAPLUS

CN 3-Pyridinecarboxamide, 6-chloro-N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-, rel- (CA INDEX NAME)

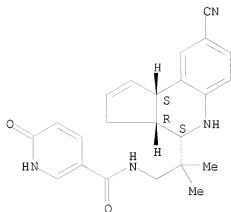
Relative stereochemistry.



RN 637333-92-7 CAPLUS

CN 3-Pyridinecarboxamide, N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-1,6-dihydro-6-oxo-, rel- (CA INDEX NAME)

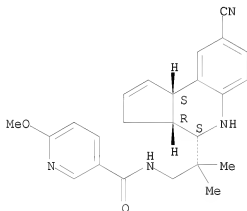
Relative stereochemistry.



RN 637333-93-8 CAPLUS

CN 3-Pyridinecarboxamide, N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-6-methoxy-, rel- (CA INDEX NAME)

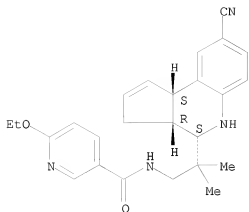
Relative stereochemistry.



RN 637333-94-9 CAPLUS

CN 3-Pyridinecarboxamide, N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-6-ethoxy-, rel- (CA INDEX NAME)

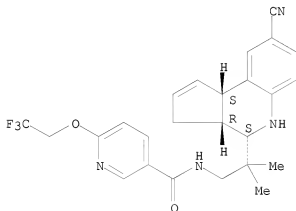
Relative stereochemistry.



RN 637333-95-0 CAPLUS

CN 3-Pyridinecarboxamide, N-[2-[(3aR, 4S, 9bS)-8-cyano-3a, 4, 5, 9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-6-(2, 2, 2-trifluoroethoxy)-, rel- (CA INDEX NAME)

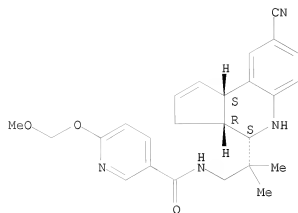
Relative stereochemistry.



RN 637333-96-1 CAPLUS

CN 3-Pyridinecarboxamide, N-[2-[(3aR, 4S, 9bS)-8-cyano-3a, 4, 5, 9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-6-(methoxymethoxy)-, rel- (CA INDEX NAME)

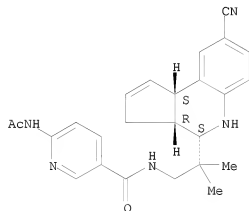
Relative stereochemistry.



RN 637333-97-2 CAPLUS

CN 3-Pyridinecarboxamide, 6-(acetylamino)-N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-, rel- (CA INDEX NAME)

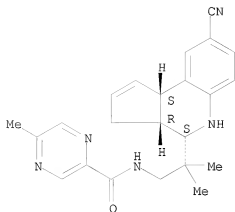
Relative stereochemistry.



RN 637333-98-3 CAPLUS

CN 2-Pyrazinecarboxamide, N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-5-methyl-, rel- (CA INDEX NAME)

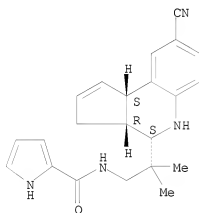
Relative stereochemistry.



RN 637333-99-4 CAPLUS

CN 1H-Pyrrole-2-carboxamide, N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-, rel- (CA INDEX NAME)

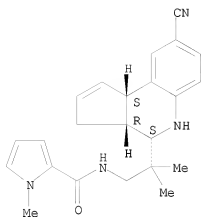
Relative stereochemistry.



RN 637334-00-0 CAPLUS

CN 1H-Pyrrole-2-carboxamide, N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-1-methyl-, rel- (CA INDEX NAME)

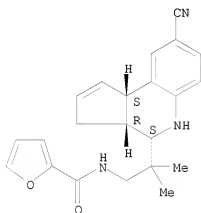
Relative stereochemistry.



RN 637334-01-1 CAPLUS

CN 2-Furancarboxamide, N-[2-[(3aR, 4S, 9bS)-8-cyano-3a, 4, 5, 9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-, rel- (CA INDEX NAME)

Relative stereochemistry.



REFERENCE COUNT: 10 THERE ARE 10 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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LOGOFF? (Y)/N/HOLD:y

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